

CERTIFICATION

| | | | |
|---------|----------|-------------|----------------------|
| SDG No: | 1701428B | Laboratory: | Eurofins, Folsom, CA |
| Site: | BMSMC | Matrix: | Air |

SUMMARY: Air samples (Table 1) were collected on the BMSMC facility. The BMSMC facility is located in Humacao, PR. Samples were taken January 24 & 26, 2017 and were analyzed in Eurofins Laboratory of Folsom, California that reported the data under SDG No.: 1701428B. Results were validated using the validation guidelines of Compendium Method TO-15. Determination of Volatile Organic Compounds (VOCs) In Air Collected In Specially-Prepared Canisters and Analyzed By Gas Chromatography/Mass Spectrometry (GC/MS), January, 1999. USEPA Hazardous Waste Support Branch. Validating Air Samples. Volatile Organic Analysis of Ambient Air in Canisters by Method TO-15, (SOP # HW-31. Revision #6. June, 2014). The analyses performed are shown in Table 1. Individual data review worksheets are enclosed for each target analyte group. The data sample summary form shows analyte results that were qualified.

In summary, the results are valid and can be used for decision making purposes.

Table 1. Samples analyzed and analysis performed

| SAMPLE ID | SAMPLE DESCRIPTION | MATRIX | ANALYSIS PERFORMED |
|--------------|--------------------|--------|--------------------|
| 1701428B-16A | B8SS-2-012417 | Air | TO-15 (full suite) |
| 1701428B-17A | B8SS-2D-012417 | Air | TO-15 (full suite) |
| 1701428B-18A | B18SS-1-012617 | Air | TO-15 (full suite) |
| 1701428B-19A | B18SS-1Dup-012617 | Air | TO-15 (full suite) |

Reviewer Name: Rafael Infante
Chemist License 1888

Signature:

Date:

March 18, 2017





Air Toxics

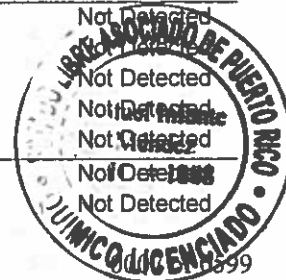
Client Sample ID: B8SS-2-012417

Lab ID#: 1701428B-16A

EPA METHOD TO-15 GC/MS

| | | |
|--------------|----------|--|
| File Name: | 14020610 | Date of Collection: 1/24/17 3:35:00 PM |
| Dil. Factor: | 24.9 | Date of Analysis: 2/6/17 02:24 PM |

| Compound | Rpt. Limit (ppbv) | Amount (ppbv) | Rpt. Limit (ug/m3) | Amount (ug/m3) |
|----------------------------------|-------------------|---------------|--------------------|----------------|
| Freon 12 | 120 | Not Detected | 620 | Not Detected |
| Freon 114 | 120 | Not Detected | 870 | Not Detected |
| Chloromethane | 500 | Not Detected | 1000 | Not Detected |
| Vinyl Chloride | 120 | Not Detected | 320 | Not Detected |
| 1,3-Butadiene | 120 | Not Detected | 280 | Not Detected |
| Bromomethane | 500 | Not Detected | 1900 | Not Detected |
| Chloroethane | 500 | Not Detected | 1300 | Not Detected |
| Freon 11 | 120 | Not Detected | 700 | Not Detected |
| Ethanol | 500 | Not Detected | 940 | Not Detected |
| Freon 113 | 120 | Not Detected | 950 | Not Detected |
| 1,1-Dichloroethene | 120 | Not Detected | 490 | Not Detected |
| Acetone | 500 | Not Detected | 1200 | Not Detected |
| 2-Propanol | 500 | Not Detected | 1200 | Not Detected |
| Carbon Disulfide | 500 | Not Detected | 1600 | Not Detected |
| 3-Chloropropene | 500 | Not Detected | 1600 | Not Detected |
| Methylene Chloride | 500 | Not Detected | 1700 | Not Detected |
| Methyl tert-butyl ether | 120 | 250 | 450 | 920 |
| trans-1,2-Dichloroethene | 120 | Not Detected | 490 | Not Detected |
| Hexane | 120 | 360 | 440 | 1300 |
| 1,1-Dichloroethane | 120 | Not Detected | 500 | Not Detected |
| 2-Butanone (Methyl Ethyl Ketone) | 500 | Not Detected | 1500 | Not Detected |
| cis-1,2-Dichloroethene | 120 | Not Detected | 490 | Not Detected |
| Tetrahydrofuran | 120 | Not Detected | 370 | Not Detected |
| Chloroform | 120 | Not Detected | 610 | Not Detected |
| 1,1,1-Trichloroethane | 120 | Not Detected | 680 | Not Detected |
| Cyclohexane | 120 | 1000 | 430 | 3600 |
| Carbon Tetrachloride | 120 | Not Detected | 780 | Not Detected |
| 2,2,4-Trimethylpentane | 120 | Not Detected | 580 | Not Detected |
| Benzene | 120 | Not Detected | 400 | Not Detected |
| 1,2-Dichloroethane | 120 | Not Detected | 500 | Not Detected |
| Heptane | 120 | 480 | 510 | 2000 |
| Trichloroethene | 120 | Not Detected | 670 | Not Detected |
| 1,2-Dichloropropane | 120 | Not Detected | 580 | Not Detected |
| 1,4-Dioxane | 500 | Not Detected | 1800 | Not Detected |
| Bromodichloromethane | 120 | Not Detected | 830 | Not Detected |
| cis-1,3-Dichloropropene | 120 | Not Detected | 560 | Not Detected |
| 4-Methyl-2-pentanone | 120 | Not Detected | 510 | Not Detected |
| Toluene | 120 | Not Detected | 470 | Not Detected |
| trans-1,3-Dichloropropene | 120 | Not Detected | 560 | Not Detected |
| 1,1,2-Trichloroethane | 120 | Not Detected | 680 | Not Detected |
| Tetrachloroethene | 120 | Not Detected | 840 | Not Detected |
| 2-Hexanone | 500 | Not Detected | 2000 | Not Detected |





Air Toxics

Client Sample ID: B8SS-2-012417

Lab ID#: 1701428B-16A

EPA METHOD TO-15 GC/MS

| | | |
|--------------|----------|--|
| File Name: | 14020610 | Date of Collection: 1/24/17 3:35:00 PM |
| Dil. Factor: | 24.9 | Date of Analysis: 2/6/17 02:24 PM |

| Compound | Rpt. Limit (ppbv) | Amount (ppbv) | Rpt. Limit (ug/m3) | Amount (ug/m3) |
|---------------------------|-------------------|---------------|--------------------|----------------|
| Dibromochloromethane | 120 | Not Detected | 1100 | Not Detected |
| 1,2-Dibromoethane (EDB) | 120 | Not Detected | 960 | Not Detected |
| Chlorobenzene | 120 | Not Detected | 570 | Not Detected |
| Ethyl Benzene | 120 | Not Detected | 540 | Not Detected |
| m,p-Xylene | 120 | 52 J | 540 | 220 J |
| o-Xylene | 120 | Not Detected | 540 | Not Detected |
| Styrene | 120 | Not Detected | 530 | Not Detected |
| Bromoform | 120 | Not Detected | 1300 | Not Detected |
| Cumene | 120 | 130 | 610 | 640 |
| 1,1,2,2-Tetrachloroethane | 120 | Not Detected | 850 | Not Detected |
| Propylbenzene | 120 | Not Detected | 610 | Not Detected |
| 4-Ethyltoluene | 120 | Not Detected | 610 | Not Detected |
| 1,3,5-Trimethylbenzene | 120 | Not Detected | 610 | Not Detected |
| 1,2,4-Trimethylbenzene | 120 | Not Detected | 610 | Not Detected |
| 1,3-Dichlorobenzene | 120 | Not Detected | 750 | Not Detected |
| 1,4-Dichlorobenzene | 120 | Not Detected | 750 | Not Detected |
| alpha-Chlorotoluene | 120 | Not Detected | 640 | Not Detected |
| 1,2-Dichlorobenzene | 120 | Not Detected | 750 | Not Detected |
| 1,2,4-Trichlorobenzene | 500 | Not Detected | 3700 | Not Detected |
| Hexachlorobutadiene | 500 | Not Detected | 5300 | Not Detected |
| Naphthalene | 500 | Not Detected | 2600 | Not Detected |

J = Estimated value.

Container Type: 1 Liter Summa Canister (100% Certified)

| Surrogates | %Recovery | Method Limits |
|-----------------------|-----------|---------------|
| 1,2-Dichloroethane-d4 | 96 | 70-130 |
| Toluene-d8 | 97 | 70-130 |
| 4-Bromofluorobenzene | 102 | 70-130 |





Air Toxics

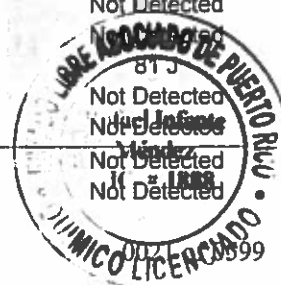
Client Sample ID: B8SS-2D-012417

Lab ID#: 1701428B-17A

EPA METHOD TO-15 GC/MS

File Name: 14020611 Date of Collection: 1/24/17 3:35:00 PM
Dil. Factor: 24.2 Date of Analysis: 2/6/17 03:00 PM

| Compound | Rpt. Limit (ppbv) | Amount (ppbv) | Rpt. Limit (ug/m3) | Amount (ug/m3) |
|----------------------------------|----------------------|------------------|-----------------------|-------------------|
| Freon 12 | 120 | Not Detected | 600 | Not Detected |
| Freon 114 | 120 | Not Detected | 840 | Not Detected |
| Chloromethane | 480 | Not Detected | 1000 | Not Detected |
| Vinyl Chloride | 120 | Not Detected | 310 | Not Detected |
| 1,3-Butadiene | 120 | Not Detected | 270 | Not Detected |
| Bromomethane | 480 | Not Detected | 1900 | Not Detected |
| Chloroethane | 480 | Not Detected | 1300 | Not Detected |
| Freon 11 | 120 | Not Detected | 680 | Not Detected |
| Ethanol | 480 | Not Detected | 910 | Not Detected |
| Freon 113 | 120 | Not Detected | 930 | Not Detected |
| 1,1-Dichloroethene | 120 | Not Detected | 480 | Not Detected |
| Acetone | 480 | Not Detected | 1100 | Not Detected |
| 2-Propanol | 480 | Not Detected | 1200 | Not Detected |
| Carbon Disulfide | 480 | Not Detected | 1500 | Not Detected |
| 3-Chloropropene | 480 | Not Detected | 1500 | Not Detected |
| Methylene Chloride | 480 | Not Detected | 1700 | Not Detected |
| Methyl tert-butyl ether | 120 | 250 | 440 | 900 |
| trans-1,2-Dichloroethene | 120 | Not Detected | 480 | Not Detected |
| Hexane | 120 | 360 | 430 | 1300 |
| 1,1-Dichloroethane | 120 | Not Detected | 490 | Not Detected |
| 2-Butanone (Methyl Ethyl Ketone) | 480 | Not Detected | 1400 | Not Detected |
| cis-1,2-Dichloroethene | 120 | Not Detected | 480 | Not Detected |
| Tetrahydrofuran | 120 | Not Detected | 360 | Not Detected |
| Chloroform | 120 | Not Detected | 590 | Not Detected |
| 1,1,1-Trichloroethane | 120 | Not Detected | 660 | Not Detected |
| Cyclohexane | 120 | 980 | 420 | 3400 |
| Carbon Tetrachloride | 120 | Not Detected | 760 | Not Detected |
| 2,2,4-Trimethylpentane | 120 | 100 J | 560 | 470 J |
| Benzene | 120 | Not Detected | 390 | Not Detected |
| 1,2-Dichloroethane | 120 | Not Detected | 490 | Not Detected |
| Heptane | 120 | 440 | 500 | 1800 |
| Trichloroethene | 120 | Not Detected | 650 | Not Detected |
| 1,2-Dichloropropane | 120 | Not Detected | 560 | Not Detected |
| 1,4-Dioxane | 480 | Not Detected | 1700 | Not Detected |
| Bromodichloromethane | 120 | Not Detected | 810 | Not Detected |
| cis-1,3-Dichloropropene | 120 | Not Detected | 550 | Not Detected |
| 4-Methyl-2-pentanone | 120 | Not Detected | 500 | Not Detected |
| Toluene | 120 | 21 J | 460 | 81 J |
| trans-1,3-Dichloropropene | 120 | Not Detected | 550 | Not Detected |
| 1,1,2-Trichloroethane | 120 | Not Detected | 660 | Not Detected |
| Tetrachloroethene | 120 | Not Detected | 820 | Not Detected |
| 2-Hexanone | 480 | Not Detected | 2000 | Not Detected |





Air Toxics

Client Sample ID: B8SS-2D-012417

Lab ID#: 1701428B-17A

EPA METHOD TO-15 GC/MS

| | | |
|--------------|----------|--|
| File Name: | 14020611 | Date of Collection: 1/24/17 3:35:00 PM |
| Dil. Factor: | 24.2 | Date of Analysis: 2/6/17 03:00 PM |

| Compound | Rpt. Limit (ppbv) | Amount (ppbv) | Rpt. Limit (ug/m3) | Amount (ug/m3) |
|---------------------------|----------------------|------------------|-----------------------|-------------------|
| Dibromochloromethane | 120 | Not Detected | 1000 | Not Detected |
| 1,2-Dibromoethane (EDB) | 120 | Not Detected | 930 | Not Detected |
| Chlorobenzene | 120 | Not Detected | 560 | Not Detected |
| Ethyl Benzene | 120 | Not Detected | 520 | Not Detected |
| m,p-Xylene | 120 | 64 J | 520 | 280 J |
| o-Xylene | 120 | Not Detected | 520 | Not Detected |
| Styrene | 120 | Not Detected | 520 | Not Detected |
| Bromoform | 120 | Not Detected | 1200 | Not Detected |
| Cumene | 120 | 120 J | 590 | 580 J |
| 1,1,2,2-Tetrachloroethane | 120 | Not Detected | 830 | Not Detected |
| Propylbenzene | 120 | Not Detected | 590 | Not Detected |
| 4-Ethyltoluene | 120 | Not Detected | 590 | Not Detected |
| 1,3,5-Trimethylbenzene | 120 | Not Detected | 590 | Not Detected |
| 1,2,4-Trimethylbenzene | 120 | Not Detected | 590 | Not Detected |
| 1,3-Dichlorobenzene | 120 | Not Detected | 730 | Not Detected |
| 1,4-Dichlorobenzene | 120 | Not Detected | 730 | Not Detected |
| alpha-Chlorotoluene | 120 | Not Detected | 630 | Not Detected |
| 1,2-Dichlorobenzene | 120 | Not Detected | 730 | Not Detected |
| 1,2,4-Trichlorobenzene | 480 | Not Detected | 3600 | Not Detected |
| Hexachlorobutadiene | 480 | Not Detected | 5200 | Not Detected |
| Naphthalene | 480 | Not Detected | 2500 | Not Detected |

J = Estimated value.

Container Type: 1 Liter Summa Canister (100% Certified)

| Surrogates | %Recovery | Method Limits |
|-----------------------|-----------|------------------|
| 1,2-Dichloroethane-d4 | 95 | 70-130 |
| Toluene-d8 | 99 | 70-130 |
| 4-Bromofluorobenzene | 100 | 70-130 |





Air Toxics

Client Sample ID: B18SS-1-012617

Lab ID#: 1701428B-18A

EPA METHOD TO-15 GC/MS FULL SCAN

File Name: a020224 Date of Collection: 1/26/17 5:30:00 PM
DIL. Factor: 2.50 Date of Analysis: 2/3/17 01:54 AM

| Compound | Rpt. Limit (ppbv) | Amount (ppbv) | Rpt. Limit (ug/m3) | Amount (ug/m3) |
|----------------------------------|-------------------|---------------|--------------------|----------------|
| Freon 12 | 1.2 | 0.63 J | 6.2 | 3.1 J |
| Freon 114 | 1.2 | Not Detected | 8.7 | Not Detected |
| Chloromethane | 12 | Not Detected | 26 | Not Detected |
| Vinyl Chloride | 1.2 | Not Detected | 3.2 | Not Detected |
| 1,3-Butadiene | 1.2 | Not Detected | 2.8 | Not Detected |
| Bromomethane | 12 | Not Detected | 48 | Not Detected |
| Chloroethane | 5.0 | Not Detected | 13 | Not Detected |
| Freon 11 | 1.2 | 1.4 | 7.0 | 7.6 |
| Ethanol | 5.0 | 140 | 9.4 | 270 |
| Freon 113 | 1.2 | Not Detected | 9.6 | Not Detected |
| 1,1-Dichloroethene | 1.2 | Not Detected | 5.0 | Not Detected |
| Acetone | 12 | 12 J | 30 | 29 J |
| 2-Propanol | 5.0 | 180 | 12 | 460 |
| Carbon Disulfide | 5.0 | 1.2 J | 16 | 3.7 J |
| 3-Chloropropene | 5.0 | Not Detected | 16 | Not Detected |
| Methylene Chloride | 12 | Not Detected | 43 | Not Detected |
| Methyl tert-butyl ether | 5.0 | Not Detected | 18 | Not Detected |
| trans-1,2-Dichloroethene | 1.2 | Not Detected | 5.0 | Not Detected |
| Hexane | 1.2 | Not Detected | 4.4 | Not Detected |
| 1,1-Dichloroethane | 1.2 | Not Detected | 5.0 | Not Detected |
| 2-Butanone (Methyl Ethyl Ketone) | 5.0 | 2.1 J | 15 | 6.2 J |
| cis-1,2-Dichloroethene | 1.2 | Not Detected | 5.0 | Not Detected |
| Tetrahydrofuran | 1.2 | Not Detected | 3.7 | Not Detected |
| Chloroform | 1.2 | Not Detected | 6.1 | Not Detected |
| 1,1,1-Trichloroethane | 1.2 | Not Detected | 6.8 | Not Detected |
| Cyclohexane | 1.2 | Not Detected | 4.3 | Not Detected |
| Carbon Tetrachloride | 1.2 | Not Detected | 7.9 | Not Detected |
| 2,2,4-Trimethylpentane | 1.2 | Not Detected | 5.8 | Not Detected |
| Benzene | 1.2 | 0.22 J | 4.0 | 0.71 J |
| 1,2-Dichloroethane | 1.2 | Not Detected | 5.0 | Not Detected |
| Heptane | 1.2 | Not Detected | 5.1 | Not Detected |
| Trichloroethene | 1.2 | Not Detected | 6.7 | Not Detected |
| 1,2-Dichloropropane | 1.2 | Not Detected | 5.8 | Not Detected |
| 1,4-Dioxane | 5.0 | Not Detected | 18 | Not Detected |
| Bromodichloromethane | 1.2 | Not Detected | 8.4 | Not Detected |
| cis-1,3-Dichloropropene | 1.2 | Not Detected | 5.7 | Not Detected |
| 4-Methyl-2-pentanone | 1.2 | Not Detected | 5.1 | Not Detected |
| Toluene | 1.2 | 1.4 | 4.7 | Not Detected |
| trans-1,3-Dichloropropene | 1.2 | Not Detected | 5.7 | Not Detected |
| 1,1,2-Trichloroethane | 1.2 | Not Detected | 6.8 | Not Detected |
| Tetrachloroethene | 1.2 | Not Detected | 8.5 | Not Detected |
| 2-Hexanone | 5.0 | Not Detected | 20 | Not Detected |





Air Toxics

Client Sample ID: B18SS-1-012617

Lab ID#: 1701428B-18A

EPA METHOD TO-15 GC/MS FULL SCAN

File Name: a020224 Date of Collection: 1/26/17 5:30:00 PM
Dil. Factor: 2.50 Date of Analysis: 2/3/17 01:54 AM

| Compound | Rpt. Limit (ppbv) | Amount (ppbv) | Rpt. Limit (ug/m3) | Amount (ug/m3) |
|---------------------------|-------------------|---------------|--------------------|----------------|
| Dibromochloromethane | 1.2 | Not Detected | 11 | Not Detected |
| 1,2-Dibromoethane (EDB) | 1.2 | Not Detected | 9.6 | Not Detected |
| Chlorobenzene | 1.2 | Not Detected | 5.8 | Not Detected |
| Ethyl Benzene | 1.2 | 0.32 J | 5.4 | 1.4 J |
| m,p-Xylene | 1.2 | 1.0 J | 5.4 | 4.6 J |
| o-Xylene | 1.2 | 0.50 J | 5.4 | 2.2 J |
| Styrene | 1.2 | Not Detected | 5.3 | Not Detected |
| Bromoform | 1.2 | Not Detected | 13 | Not Detected |
| Cumene | 1.2 | Not Detected | 6.1 | Not Detected |
| 1,1,2,2-Tetrachloroethane | 1.2 | Not Detected | 8.6 | Not Detected |
| Propylbenzene | 1.2 | Not Detected | 6.1 | Not Detected |
| 4-Ethyltoluene | 1.2 | 0.23 J | 6.1 | 1.1 J |
| 1,3,5-Trimethylbenzene | 1.2 | Not Detected | 6.1 | Not Detected |
| 1,2,4-Trimethylbenzene | 1.2 | 0.25 J | 6.1 | 1.2 J |
| 1,3-Dichlorobenzene | 1.2 | 1.2 J | 7.5 | 7.2 J |
| 1,4-Dichlorobenzene | 1.2 | Not Detected | 7.5 | Not Detected |
| alpha-Chlorotoluene | 1.2 | Not Detected | 6.5 | Not Detected |
| 1,2-Dichlorobenzene | 1.2 | Not Detected | 7.5 | Not Detected |
| 1,2,4-Trichlorobenzene | 5.0 | Not Detected | 37 | Not Detected |
| Hexachlorobutadiene | 5.0 | Not Detected | 53 | Not Detected |
| Naphthalene | 2.5 | Not Detected | 13 | Not Detected |

J = Estimated value.

Container Type: 1 Liter Summa Canister (100% Certified)

| Surrogates | %Recovery | Method Limits |
|-----------------------|-----------|---------------|
| Toluene-d8 | 100 | 70-130 |
| 1,2-Dichloroethane-d4 | 98 | 70-130 |
| 4-Bromofluorobenzene | 106 | 70-130 |





Lab ID#: 1701428B-19A

EPA METHOD TO-15 GC/MS FULL SCAN

| | | | | |
|----------------------------------|----------------------|---------------------|-----------------------|-------------------|
| File Name: | a020311 | Date of Collection: | 1/26/17 5:30:00 PM | |
| Dil. Factor: | 2.46 | Date of Analysis: | 2/3/17 06:37 PM | |
| Compound | Rpt. Limit (ppbv) | Amount (ppbv) | Rpt. Limit (ug/m3) | Amount (ug/m3) |
| Freon 12 | 1.2 | 0.50 J | 6.1 | 2.5 J |
| Freon 114 | 1.2 | Not Detected | 8.6 | Not Detected |
| Chloromethane | 12 | Not Detected | 25 | Not Detected |
| Vinyl Chloride | 1.2 | Not Detected | 3.1 | Not Detected |
| 1,3-Butadiene | 1.2 | Not Detected | 2.7 | Not Detected |
| Bromomethane | 12 | Not Detected | 48 | Not Detected |
| Chloroethane | 4.9 | Not Detected | 13 | Not Detected |
| Freon 11 | 1.2 | 1.6 | 6.9 | 9.2 |
| Ethanol | 4.9 | 240 | 9.3 | 460 |
| Freon 113 | 1.2 | Not Detected | 9.4 | Not Detected |
| 1,1-Dichloroethene | 1.2 | Not Detected | 4.9 | Not Detected |
| Acetone | 12 | 20 | 29 | 47 |
| 2-Propanol | 4.9 | 360 | 12 | 880 |
| Carbon Disulfide | 4.9 | Not Detected | 15 | Not Detected |
| 3-Chloropropene | 4.9 | Not Detected | 15 | Not Detected |
| Methylene Chloride | 12 | Not Detected | 43 | Not Detected |
| Methyl tert-butyl ether | 4.9 | Not Detected | 18 | Not Detected |
| trans-1,2-Dichloroethene | 1.2 | Not Detected | 4.9 | Not Detected |
| Hexane | 1.2 | 0.37 J | 4.3 | 1.3 J |
| 1,1-Dichloroethane | 1.2 | Not Detected | 5.0 | Not Detected |
| 2-Butanone (Methyl Ethyl Ketone) | 4.9 | 4.6 J | 14 | 13 J |
| cis-1,2-Dichloroethene | 1.2 | Not Detected | 4.9 | Not Detected |
| Tetrahydrofuran | 1.2 | 0.15 J | 3.6 | 0.45 J |
| Chloroform | 1.2 | Not Detected | 6.0 | Not Detected |
| 1,1,1-Trichloroethane | 1.2 | Not Detected | 6.7 | Not Detected |
| Cyclohexane | 1.2 | Not Detected | 4.2 | Not Detected |
| Carbon Tetrachloride | 1.2 | Not Detected | 7.7 | Not Detected |
| 2,2,4-Trimethylpentane | 1.2 | 0.32 J | 5.7 | 1.5 J |
| Benzene | 1.2 | 0.21 J | 3.9 | 0.67 J |
| 1,2-Dichloroethane | 1.2 | Not Detected | 5.0 | Not Detected |
| Heptane | 1.2 | 0.48 J | 5.0 | 2.0 J |
| Trichloroethene | 1.2 | Not Detected | 6.6 | Not Detected |
| 1,2-Dichloropropane | 1.2 | Not Detected | 5.7 | Not Detected |
| 1,4-Dioxane | 4.9 | Not Detected | 18 | Not Detected |
| Bromodichloromethane | 1.2 | Not Detected | 8.2 | Not Detected |
| cis-1,3-Dichloropropene | 1.2 | Not Detected | 5.6 | Not Detected |
| 4-Methyl-2-pentanone | 1.2 | 1.9 | 5.0 | 7.8 |
| Toluene | 1.2 | 2.2 | 4.6 | 8.1 |
| trans-1,3-Dichloropropene | 1.2 | 3.7 | 5.7 | 17 |
| 1,1,2-Trichloroethane | 1.2 | 0.58 J | 6.7 | 1 J |
| Tetrachloroethene | 1.2 | Not Detected | 8.3 | Not Detected |
| 2-Hexanone | 4.9 | 0.54 J | 20 | 2 J |



Air Toxics

Client Sample ID: B18SS-1Dup-012617

Lab ID#: 1701428B-19A

EPA METHOD TO-15 GC/MS FULL SCAN

File Name: a020311 Date of Collection: 1/26/17 5:30:00 PM
Dil. Factor: 2.46 Date of Analysis: 2/3/17 06:37 PM

| Compound | Rpt. Limit (ppbv) | Amount (ppbv) | Rpt. Limit (ug/m3) | Amount (ug/m3) |
|---------------------------|-------------------|---------------|--------------------|----------------|
| Dibromochloromethane | 1.2 | Not Detected | 10 | Not Detected |
| 1,2-Dibromoethane (EDB) | 1.2 | Not Detected | 9.4 | Not Detected |
| Chlorobenzene | 1.2 | Not Detected | 5.7 | Not Detected |
| Ethyl Benzene | 1.2 | 0.47 J | 5.3 | 2.0 J |
| m,p-Xylene | 1.2 | 1.8 | 5.3 | 7.8 |
| o-Xylene | 1.2 | 0.89 J | 5.3 | 3.9 J |
| Styrene | 1.2 | 0.15 J | 5.2 | 0.64 J |
| Bromoform | 1.2 | Not Detected | 13 | Not Detected |
| Cumene | 1.2 | 0.29 J | 6.0 | 1.4 J |
| 1,1,2,2-Tetrachloroethane | 1.2 | 0.070 J | 8.4 | 0.48 J |
| Propylbenzene | 1.2 | 0.13 J | 6.0 | 0.66 J |
| 4-Ethyltoluene | 1.2 | 0.62 J | 6.0 | 3.1 J |
| 1,3,5-Trimethylbenzene | 1.2 | 0.15 J | 6.0 | 0.74 J |
| 1,2,4-Trimethylbenzene | 1.2 | 0.55 J | 6.0 | 2.7 J |
| 1,3-Dichlorobenzene | 1.2 | 2.5 | 7.4 | 15 |
| 1,4-Dichlorobenzene | 1.2 | 2.4 | 7.4 | 14 |
| alpha-Chlorotoluene | 1.2 | Not Detected | 6.4 | Not Detected |
| 1,2-Dichlorobenzene | 1.2 | Not Detected | 7.4 | Not Detected |
| 1,2,4-Trichlorobenzene | 4.9 | Not Detected | 36 | Not Detected |
| Hexachlorobutadiene | 4.9 | Not Detected | 52 | Not Detected |
| Naphthalene | 2.5 | Not Detected | 13 | Not Detected |

J = Estimated value.

Container Type: 1 Liter Summa Canister (100% Certified)

| Surrogates | %Recovery | Method Limits |
|-----------------------|-----------|---------------|
| Toluene-d8 | 102 | 70-130 |
| 1,2-Dichloroethane-d4 | 100 | 70-130 |
| 4-Bromofluorobenzene | 102 | 70-130 |





Chain-of-Custody Record

Page 2 of 2

Contact Person Terry Taylor
Company Anderson Mulholland & Associates, Inc.
Address 2700 Winchester, Suite 41 City Purchase State NY Zip 10577
Phone 914-251-0400, x 309 FAX
Collected By: (Signature) *Terry Taylor*

Project Information:
P.O. #
Project # BMS VI Invest.
Project Name
Buildings 8, 13, 5, 18, and 30

Turn Around Time:
☒ Normal
☐ Rush
Specify

Pressurized
by:
Date:
Press. Gas: N2 He

| Lab I.D. | Field Sample I.D. | Canister I.D. | Date & Time | Analysis Requested | Canister Initial | Canister Final | Pressure/Vacuum Receipt | Final (psi) |
|----------|-------------------|---------------|--------------|------------------------------|------------------|----------------|-------------------------|-------------|
| 117 | B181A-3-012117 | 34320 | 1/21/13:1510 | TO-15, MeOH, CH ₄ | 29 "Hg | 6 | | |
| 118 | B181A-4-012117 | U1591 | 1/21/13:1512 | TO-15, MeOH, CH ₄ | 29.5 "Hg | 6 | | |
| 119 | B81A-2-012317 | U1726 | 1/23/17:1223 | TO-15, MeOH, CH ₄ | 29 "Hg | 8 | | |
| 120 | B81A-2D-012317 | U6416 | 1/23/17:1823 | TO-15, MeOH, CH ₄ | 29 "Hg | 9 | | |
| 121 | B8AA-012317 | 00316 | 1/23/17:0920 | TO-15, MeOH, CH ₄ | 29 "Hg | 3 | | |
| 122 | B855-2-012417 | 1L1603 | 1/24/17:1535 | TO-15, MeOH, CH ₄ | 30 "Hg | 5 | | |
| 123 | B855-2D-012417 | 12323 | 1/24/17:1535 | TO-15, MeOH, CH ₄ | 29 "Hg | 5 | | |
| 124 | B1855-1-012617 | 1L1720 | 1/26/17:1930 | TO-15, MeOH, CH ₄ | 30 "Hg | 5 | | |
| 125 | B1855-1000-012617 | | | | "Hg | | | |
| 126 | B1855-1000-012617 | U1929 | 1/26/17:1930 | TO-15, MeOH, CH ₄ | 30 "Hg | 6 | | |

Relinquished By: (Signature) *Terry Taylor* Date/Time 1-27-17 1400
Relinquished By: (Signature) Date/Time
Relinquished By: (Signature) Date/Time
Relinquished By: (Signature) Date/Time

Notes:
FedEx Tracking No. 7782 2053 0772

Shipper Name Air Bill # Opened By Temp @ Condition Custody Seals Work Order #
Fed Ex 5772 N/A good Yes No None 1701428

EXECUTIVE NARRATIVE

SDG No: **1701428B** Laboratory: **Eurofins, Folsom, CA**
Analysis: **TO-15** Number of Samples: **4**
Location:

SUMMARY: Four (4) samples were analyzed for VOCs in ambient air following Compendium Method TO-15. The sample results were assessed according to USEPA data validation guidance documents in the following order of precedence: "Compendium Method TO-15. Determination of Volatile Organic Compounds (VOCs) In Air Collected In Specially-Prepared Canisters and Analyzed By Gas Chromatography/Mass Spectrometry (GC/MS), January, 1999"; USEPA Hazardous Waste Support Branch. Validating Air Samples. Volatile Organic Analysis of Ambient Air in Canisters by Method TO-15, (SOP # HW-31. Revision #6. June, 2014). The QC criteria and data validation actions listed on the data review worksheets are from the primary guidance document, unless otherwise noted.

Results are valid and can be used for decision making purposes.

Critical issues: **None**
Major: **None**
Minor: **None**

Critical findings: **None**
Major findings: **None**
Minor findings:

1. Initial and continuing calibrations meet method specific requirements except for the cases described in the Data Review Worksheet. For analytes not meeting the % difference continuing calibration criteria their results qualified as estimated (J or UJ) in affected samples.
2. Several analytes detected at the method blank analyzed on 02/03/17.
3. LCS/LCSD (Blank spike) analyzed in this data package. % recoveries and RPD within laboratory control limits except for the cases described in the Data Review Worksheet. Positive results for Freon 113 and Naphthalene qualified as estimated (J).
4. Laboratory/field duplicate analyzed with this data package. RPD within laboratory and method performed criteria for analytes with concentration > 5 x SQL except for the cases described in the Data Review Worksheet. Results for acetone and benzene qualified (U) in affected samples as per Table 7. Blank actions for TO-15 Analyses.
5. Results for ethanol and 2-propanol qualified as estimated (J) in samples 1701428B-18A/1701428B-19A due to RPD exceeding the method performance criteria.

COMMENTS: Results are valid and can be used for decision making purposes.

Reviewers Name: Rafael Infante
Chemist License 1888



Signature:

Date: March 15, 2017

TO-15 DATA SAMPLE SUMMARY

Sample ID: 1701428B-16A

Sample location: BMSMC, Humacao, PR

Sampling date: 24-Jan-17

Matrix: Air

METHOD: TO-15

| Analyte Name | Result | Units | Dilution Factor | Lab Flag | Validation | Reportable |
|----------------------------------|--------|-------------------|-----------------|----------|------------|------------|
| Freon 12 | 620 | ug/m ³ | 24.9 | - | U | Yes |
| Freon 114 | 870 | ug/m ³ | 24.9 | - | U | Yes |
| Chloromethane | 1000 | ug/m ³ | 24.9 | - | U | Yes |
| Vinyl Chloride | 320 | ug/m ³ | 24.9 | - | U | Yes |
| 1,3-Butadiene | 280 | ug/m ³ | 24.9 | - | U | Yes |
| Bromomethane | 1900 | ug/m ³ | 24.9 | - | U | Yes |
| Chloroethane | 1300 | ug/m ³ | 24.9 | - | U | Yes |
| Freon 11 | 700 | ug/m ³ | 24.9 | - | UJ | Yes |
| Ethanol | 940 | ug/m ³ | 24.9 | - | U | Yes |
| Freon 113 | 950 | ug/m ³ | 24.9 | - | UJ | Yes |
| 1,1-Dichloroethene | 490 | ug/m ³ | 24.9 | - | U | Yes |
| Acetone | 1200 | ug/m ³ | 24.9 | - | U | Yes |
| 2-Propanol | 1200 | ug/m ³ | 24.9 | - | U | Yes |
| Carbon Disulfide | 1600 | ug/m ³ | 24.9 | - | U | Yes |
| 3-Chloropropene | 1600 | ug/m ³ | 24.9 | - | U | Yes |
| Methylene Chloride | 1700 | ug/m ³ | 24.9 | - | U | Yes |
| Methyl tert-butyl ether | 920 | ug/m ³ | 24.9 | - | - | Yes |
| trans-1,2-Dichloroethene | 490 | ug/m ³ | 24.9 | - | U | Yes |
| Hexane | 1300 | ug/m ³ | 24.9 | - | - | Yes |
| 1,1-Dichloroethane | 500 | ug/m ³ | 24.9 | - | U | Yes |
| 2-Butanone (Methyl Ethyl Ketone) | 1500 | ug/m ³ | 24.9 | - | U | Yes |
| cis-1,2-Dichloroethene | 490 | ug/m ³ | 24.9 | - | U | Yes |
| Tetrahydrofuran | 370 | ug/m ³ | 24.9 | - | U | Yes |
| Chloroform | 610 | ug/m ³ | 24.9 | - | U | Yes |
| 1,1,1-Trichloroethane | 680 | ug/m ³ | 24.9 | - | U | Yes |
| Cyclohexane | 3600 | ug/m ³ | 24.9 | - | - | Yes |
| Carbon Tetrachloride | 780 | ug/m ³ | 24.9 | - | U | Yes |
| 2,2,4-Trimethylpentane | 580 | ug/m ³ | 24.9 | - | U | Yes |
| Benzene | 400 | ug/m ³ | 24.9 | - | U | Yes |
| 1,2-Dichloroethane | 500 | ug/m ³ | 24.9 | - | U | Yes |
| Heptane | 2000 | ug/m ³ | 24.9 | - | - | Yes |
| Trichloroethene | 670 | ug/m ³ | 24.9 | - | U | Yes |
| 1,2-Dichloropropane | 580 | ug/m ³ | 24.9 | - | U | Yes |
| 1,4-Dioxane | 1800 | ug/m ³ | 24.9 | - | U | Yes |
| Bromodichloromethane | 830 | ug/m ³ | 24.9 | - | U | Yes |
| cis-1,3-Dichloropropene | 560 | ug/m ³ | 24.9 | - | U | Yes |

| | | | | | | |
|---------------------------|------|-------------------|------|---|---|-----|
| 4-Methyl-2-pentanone | 510 | ug/m ³ | 24.9 | - | U | Yes |
| Toluene | 470 | ug/m ³ | 24.9 | - | U | Yes |
| trans-1,3-Dichloropropene | 560 | ug/m ³ | 24.9 | - | U | Yes |
| 1,1,2-Trichloroethane | 680 | ug/m ³ | 24.9 | - | U | Yes |
| Tetrachloroethene | 840 | ug/m ³ | 24.9 | - | U | Yes |
| 2-Hexanone | 2000 | ug/m ³ | 24.9 | - | U | Yes |
| Dibromochloromethane | 1100 | ug/m ³ | 24.9 | - | U | Yes |
| 1,2-Dibromoethane (EDB) | 960 | ug/m ³ | 24.9 | - | U | Yes |
| Chlorobenzene | 570 | ug/m ³ | 24.9 | - | U | Yes |
| Ethyl Benzene | 540 | ug/m ³ | 24.9 | - | U | Yes |
| m,p-Xylene | 220 | ug/m ³ | 24.9 | J | J | Yes |
| o-Xylene | 540 | ug/m ³ | 24.9 | - | U | Yes |
| Styrene | 530 | ug/m ³ | 24.9 | - | U | Yes |
| Bromoform | 1300 | ug/m ³ | 24.9 | - | U | Yes |
| Cumene | 640 | ug/m ³ | 24.9 | - | - | Yes |
| 1,1,2,2-Tetrachloroethane | 850 | ug/m ³ | 24.9 | - | U | Yes |
| Propylbenzene | 610 | ug/m ³ | 24.9 | - | U | Yes |
| 4-Ethyltoluene | 610 | ug/m ³ | 24.9 | - | U | Yes |
| 1,3,5-Trimethylbenzene | 610 | ug/m ³ | 24.9 | - | U | Yes |
| 1,2,4-Trimethylbenzene | 610 | ug/m ³ | 24.9 | - | U | Yes |
| 1,3-Dichlorobenzene | 750 | ug/m ³ | 24.9 | - | U | Yes |
| 1,4-Dichlorobenzene | 750 | ug/m ³ | 24.9 | - | U | Yes |
| alpha-Chlorotoluene | 640 | ug/m ³ | 24.9 | - | U | Yes |
| 1,2-Dichlorobenzene | 750 | ug/m ³ | 24.9 | - | U | Yes |
| 1,2,4-Trichlorobenzene | 3700 | ug/m ³ | 24.9 | - | U | Yes |
| Hexachlorobutadiene | 5300 | ug/m ³ | 24.9 | - | U | Yes |
| Naphthalene | 2600 | ug/m ³ | 24.9 | - | U | Yes |

Sample ID: 1701428B-17A
Sample location: BMSMC, Humacao, PR
Sampling date: 24-Jan-17
Matrix: Air

METHOD: TO-15

| Analyte Name | Result | Units | Dilution Factor | Lab Flag | Validation | Reportable |
|----------------------------------|--------|-------------------|-----------------|----------|------------|------------|
| Freon 12 | 600 | ug/m ³ | 24.2 | - | U | Yes |
| Freon 114 | 840 | ug/m ³ | 24.2 | - | U | Yes |
| Chloromethane | 1000 | ug/m ³ | 24.2 | - | U | Yes |
| Vinyl Chloride | 310 | ug/m ³ | 24.2 | - | U | Yes |
| 1,3-Butadiene | 270 | ug/m ³ | 24.2 | - | U | Yes |
| Bromomethane | 1900 | ug/m ³ | 24.2 | - | U | Yes |
| Chloroethane | 1300 | ug/m ³ | 24.2 | - | U | Yes |
| Freon 11 | 680 | ug/m ³ | 24.2 | - | UJ | Yes |
| Ethanol | 910 | ug/m ³ | 24.2 | - | U | Yes |
| Freon 113 | 930 | ug/m ³ | 24.2 | - | UJ | Yes |
| 1,1-Dichloroethene | 480 | ug/m ³ | 24.2 | - | U | Yes |
| Acetone | 1100 | ug/m ³ | 24.2 | - | U | Yes |
| 2-Propanol | 1200 | ug/m ³ | 24.2 | - | U | Yes |
| Carbon Disulfide | 1500 | ug/m ³ | 24.2 | - | U | Yes |
| 3-Chloropropene | 1500 | ug/m ³ | 24.2 | - | U | Yes |
| Methylene Chloride | 1700 | ug/m ³ | 24.2 | - | U | Yes |
| Methyl tert-butyl ether | 900 | ug/m ³ | 24.2 | - | - | Yes |
| trans-1,2-Dichloroethene | 480 | ug/m ³ | 24.2 | - | U | Yes |
| Hexane | 1300 | ug/m ³ | 24.2 | - | - | Yes |
| 1,1-Dichloroethane | 490 | ug/m ³ | 24.2 | - | U | Yes |
| 2-Butanone (Methyl Ethyl Ketone) | 1400 | ug/m ³ | 24.2 | - | U | Yes |
| cis-1,2-Dichloroethene | 480 | ug/m ³ | 24.2 | - | U | Yes |
| Tetrahydrofuran | 360 | ug/m ³ | 24.2 | - | U | Yes |
| Chloroform | 590 | ug/m ³ | 24.2 | - | U | Yes |
| 1,1,1-Trichloroethane | 660 | ug/m ³ | 24.2 | - | U | Yes |
| Cyclohexane | 3400 | ug/m ³ | 24.2 | - | - | Yes |
| Carbon Tetrachloride | 760 | ug/m ³ | 24.2 | - | U | Yes |
| 2,2,4-Trimethylpentane | 470 | ug/m ³ | 24.2 | J | J | Yes |
| Benzene | 390 | ug/m ³ | 24.2 | - | U | Yes |
| 1,2-Dichloroethane | 490 | ug/m ³ | 24.2 | - | U | Yes |
| Heptane | 1800 | ug/m ³ | 24.2 | - | - | Yes |
| Trichloroethene | 650 | ug/m ³ | 24.2 | - | U | Yes |
| 1,2-Dichloropropane | 560 | ug/m ³ | 24.2 | - | U | Yes |
| 1,4-Dioxane | 1700 | ug/m ³ | 24.2 | - | U | Yes |
| Bromodichloromethane | 810 | ug/m ³ | 24.2 | - | U | Yes |
| cis-1,3-Dichloropropene | 550 | ug/m ³ | 24.2 | - | U | Yes |
| 4-Methyl-2-pentanone | 500 | ug/m ³ | 24.2 | - | U | Yes |
| Toluene | 81 | ug/m ³ | 24.2 | J | J | Yes |

| | | | | | | |
|---------------------------|------|-------------------|------|---|---|-----|
| trans-1,3-Dichloropropene | 550 | ug/m ³ | 24.2 | - | U | Yes |
| 1,1,2-Trichloroethane | 660 | ug/m ³ | 24.2 | - | U | Yes |
| Tetrachloroethene | 820 | ug/m ³ | 24.2 | - | U | Yes |
| 2-Hexanone | 2000 | ug/m ³ | 24.2 | - | U | Yes |
| Dibromochloromethane | 1000 | ug/m ³ | 24.2 | - | U | Yes |
| 1,2-Dibromoethane (EDB) | 930 | ug/m ³ | 24.2 | - | U | Yes |
| Chlorobenzene | 560 | ug/m ³ | 24.2 | - | U | Yes |
| Ethyl Benzene | 520 | ug/m ³ | 24.2 | - | U | Yes |
| m,p-Xylene | 280 | ug/m ³ | 24.2 | J | J | Yes |
| o-Xylene | 520 | ug/m ³ | 24.2 | - | U | Yes |
| Styrene | 520 | ug/m ³ | 24.2 | - | U | Yes |
| Bromoform | 1200 | ug/m ³ | 24.2 | - | U | Yes |
| Cumene | 580 | ug/m ³ | 24.2 | J | J | Yes |
| 1,1,2,2-Tetrachloroethane | 830 | ug/m ³ | 24.2 | - | U | Yes |
| Propylbenzene | 590 | ug/m ³ | 24.2 | - | U | Yes |
| 4-Ethyltoluene | 590 | ug/m ³ | 24.2 | - | U | Yes |
| 1,3,5-Trimethylbenzene | 590 | ug/m ³ | 24.2 | - | U | Yes |
| 1,2,4-Trimethylbenzene | 590 | ug/m ³ | 24.2 | - | U | Yes |
| 1,3-Dichlorobenzene | 730 | ug/m ³ | 24.2 | - | U | Yes |
| 1,4-Dichlorobenzene | 730 | ug/m ³ | 24.2 | - | U | Yes |
| alpha-Chlorotoluene | 630 | ug/m ³ | 24.2 | - | U | Yes |
| 1,2-Dichlorobenzene | 730 | ug/m ³ | 24.2 | - | U | Yes |
| 1,2,4-Trichlorobenzene | 3600 | ug/m ³ | 24.2 | - | U | Yes |
| Hexachlorobutadiene | 5200 | ug/m ³ | 24.2 | - | U | Yes |
| Naphthalene | 2500 | ug/m ³ | 24.2 | - | U | Yes |

Sample ID: 1701428B-18A
Sample location: BMSMC, Humacao, PR
Sampling date: 26-Jan-17
Matrix: Air

METHOD: TO-15

| Analyte Name | Result | Units | Dilution Factor | Lab Flag | Validation | Reportable |
|----------------------------------|--------|-------------------|-----------------|----------|------------|------------|
| Freon 12 | 3.1 | ug/m ³ | 2.50 | J | J | Yes |
| Freon 114 | 8.7 | ug/m ³ | 2.50 | - | U | Yes |
| Chloromethane | 26 | ug/m ³ | 2.50 | - | U | Yes |
| Vinyl Chloride | 3.2 | ug/m ³ | 2.50 | - | U | Yes |
| 1,3-Butadiene | 2.8 | ug/m ³ | 2.50 | - | U | Yes |
| Bromomethane | 48 | ug/m ³ | 2.50 | - | U | Yes |
| Chloroethane | 13 | ug/m ³ | 2.50 | - | U | Yes |
| Freon 11 | 7.6 | ug/m ³ | 2.50 | - | - | Yes |
| Ethanol | 270 | ug/m ³ | 2.50 | - | J | Yes |
| Freon 113 | 9.6 | ug/m ³ | 2.50 | - | U | Yes |
| 1,1-Dichloroethene | 5.0 | ug/m ³ | 2.50 | - | U | Yes |
| Acetone | 30 | ug/m ³ | 2.50 | J | U | Yes |
| 2-Propanol | 460 | ug/m ³ | 2.50 | - | J | Yes |
| Carbon Disulfide | 3.7 | ug/m ³ | 2.50 | J | J | Yes |
| 3-Chloropropene | 16.0 | ug/m ³ | 2.50 | - | U | Yes |
| Methylene Chloride | 43 | ug/m ³ | 2.50 | - | U | Yes |
| Methyl tert-butyl ether | 18 | ug/m ³ | 2.50 | - | U | Yes |
| trans-1,2-Dichloroethene | 5.0 | ug/m ³ | 2.50 | - | U | Yes |
| Hexane | 4.4 | ug/m ³ | 2.50 | - | U | Yes |
| 1,1-Dichloroethane | 5.0 | ug/m ³ | 2.50 | - | U | Yes |
| 2-Butanone (Methyl Ethyl Ketone) | 6.2 | ug/m ³ | 2.50 | J | J | Yes |
| cis-1,2-Dichloroethene | 5.0 | ug/m ³ | 2.50 | - | U | Yes |
| Tetrahydrofuran | 3.7 | ug/m ³ | 2.50 | - | U | Yes |
| Chloroform | 6.10 | ug/m ³ | 2.50 | - | U | Yes |
| 1,1,1-Trichloroethane | 6.8 | ug/m ³ | 2.50 | - | U | Yes |
| Cyclohexane | 4.3 | ug/m ³ | 2.50 | - | U | Yes |
| Carbon Tetrachloride | 7.9 | ug/m ³ | 2.50 | J | J | Yes |
| 2,2,4-Trimethylpentane | 7.8 | ug/m ³ | 2.50 | - | U | Yes |
| Benzene | 4.0 | ug/m ³ | 2.50 | J | U | Yes |
| 1,2-Dichloroethane | 5.0 | ug/m ³ | 2.50 | - | U | Yes |
| Heptane | 5.10 | ug/m ³ | 2.50 | - | U | Yes |
| Trichloroethene | 6.7 | ug/m ³ | 2.50 | - | U | Yes |
| 1,2-Dichloropropane | 5.8 | ug/m ³ | 2.50 | - | U | Yes |
| 1,4-Dioxane | 18 | ug/m ³ | 2.50 | - | U | Yes |
| Bromodichloromethane | 8.4 | ug/m ³ | 2.50 | - | U | Yes |
| cis-1,3-Dichloropropene | 5.7 | ug/m ³ | 2.50 | - | U | Yes |
| 4-Methyl-2-pentanone | 5.1 | ug/m ³ | 2.50 | - | U | Yes |
| Toluene | 5.4 | ug/m ³ | 2.50 | - | - | Yes |

| | | | | | | |
|---------------------------|-----|-------------------|------|---|---|-----|
| trans-1,3-Dichloropropene | 5.7 | ug/m ³ | 2.50 | - | U | Yes |
| 1,1,2-Trichloroethane | 6.8 | ug/m ³ | 2.50 | - | U | Yes |
| Tetrachloroethene | 8.5 | ug/m ³ | 2.50 | - | U | Yes |
| 2-Hexanone | 20 | ug/m ³ | 2.50 | - | U | Yes |
| Dibromochloromethane | 11 | ug/m ³ | 2.50 | - | U | Yes |
| 1,2-Dibromoethane (EDB) | 9.6 | ug/m ³ | 2.50 | - | U | Yes |
| Chlorobenzene | 5.8 | ug/m ³ | 2.50 | - | U | Yes |
| Ethyl Benzene | 1.4 | ug/m ³ | 2.50 | J | J | Yes |
| m,p-Xylene | 4.6 | ug/m ³ | 2.50 | J | J | Yes |
| o-Xylene | 2.2 | ug/m ³ | 2.50 | J | J | Yes |
| Styrene | 5.3 | ug/m ³ | 2.50 | - | U | Yes |
| Bromoform | 13 | ug/m ³ | 2.50 | - | U | Yes |
| Cumene | 6.1 | ug/m ³ | 2.50 | - | U | Yes |
| 1,1,2,2-Tetrachloroethane | 8.6 | ug/m ³ | 2.50 | - | U | Yes |
| Propylbenzene | 6.1 | ug/m ³ | 2.50 | - | U | Yes |
| 4-Ethyltoluene | 1.1 | ug/m ³ | 2.50 | J | J | Yes |
| 1,3,5-Trimethylbenzene | 6.1 | ug/m ³ | 2.50 | - | U | Yes |
| 1,2,4-Trimethylbenzene | 1.2 | ug/m ³ | 2.50 | J | J | Yes |
| 1,3-Dichlorobenzene | 7.2 | ug/m ³ | 2.50 | J | J | Yes |
| 1,4-Dichlorobenzene | 7.5 | ug/m ³ | 2.50 | - | U | Yes |
| alpha-Chlorotoluene | 6.5 | ug/m ³ | 2.50 | - | U | Yes |
| 1,2-Dichlorobenzene | 7.5 | ug/m ³ | 2.50 | - | U | Yes |
| 1,2,4-Trichlorobenzene | 37 | ug/m ³ | 2.50 | - | U | Yes |
| Hexachlorobutadiene | 53 | ug/m ³ | 2.50 | - | U | Yes |
| Naphthalene | 13 | ug/m ³ | 2.50 | J | U | Yes |

Sample ID: 1701428B-19A
Sample location: BMSMC, Humacao, PR
Sampling date: 26-Jan-17
Matrix: Air

METHOD: TO-15

| Analyte Name | Result | Units | Dilution Factor | Lab Flag | Validation | Reportable |
|----------------------------------|--------|-------------------|-----------------|----------|------------|------------|
| Freon 12 | 2.5 | ug/m ³ | 2.46 | J | J | Yes |
| Freon 114 | 8.6 | ug/m ³ | 2.46 | - | U | Yes |
| Chloromethane | 25 | ug/m ³ | 2.46 | - | U | Yes |
| Vinyl Chloride | 3.1 | ug/m ³ | 2.46 | - | U | Yes |
| 1,3-Butadiene | 2.7 | ug/m ³ | 2.46 | - | U | Yes |
| Bromomethane | 48 | ug/m ³ | 2.46 | - | U | Yes |
| Chloroethane | 13 | ug/m ³ | 2.46 | - | U | Yes |
| Freon 11 | 9.2 | ug/m ³ | 2.46 | - | - | Yes |
| Ethanol | 460 | ug/m ³ | 2.46 | - | J | Yes |
| Freon 113 | 9.4 | ug/m ³ | 2.46 | - | U | Yes |
| 1,1-Dichloroethene | 4.9 | ug/m ³ | 2.46 | - | U | Yes |
| Acetone | 47 | ug/m ³ | 2.46 | - | U | Yes |
| 2-Propanol | 880 | ug/m ³ | 2.46 | - | J | Yes |
| Carbon Disulfide | 15 | ug/m ³ | 2.46 | - | U | Yes |
| 3-Chloropropene | 15 | ug/m ³ | 2.46 | - | U | Yes |
| Methylene Chloride | 43 | ug/m ³ | 2.46 | - | U | Yes |
| Methyl tert-butyl ether | 18 | ug/m ³ | 2.46 | - | U | Yes |
| trans-1,2-Dichloroethene | 4.9 | ug/m ³ | 2.46 | - | U | Yes |
| Hexane | 1.3 | ug/m ³ | 2.46 | J | J | Yes |
| 1,1-Dichloroethane | 5.0 | ug/m ³ | 2.46 | - | U | Yes |
| 2-Butanone (Methyl Ethyl Ketone) | 13 | ug/m ³ | 2.46 | J | J | Yes |
| cis-1,2-Dichloroethene | 4.9 | ug/m ³ | 2.46 | - | U | Yes |
| Tetrahydrofuran | 0.45 | ug/m ³ | 2.46 | J | J | Yes |
| Chloroform | 6.0 | ug/m ³ | 2.46 | - | U | Yes |
| 1,1,1-Trichloroethane | 6.7 | ug/m ³ | 2.46 | - | U | Yes |
| Cyclohexane | 4.2 | ug/m ³ | 2.46 | - | U | Yes |
| Carbon Tetrachloride | 7.7 | ug/m ³ | 2.46 | - | U | Yes |
| 2,2,4-Trimethylpentane | 1.5 | ug/m ³ | 2.46 | J | J | Yes |
| Benzene | 3.9 | ug/m ³ | 2.46 | J | U | Yes |
| 1,2-Dichloroethane | 5.0 | ug/m ³ | 2.46 | - | U | Yes |
| Heptane | 2.0 | ug/m ³ | 2.46 | J | J | Yes |
| Trichloroethene | 6.6 | ug/m ³ | 2.46 | - | U | Yes |
| 1,2-Dichloropropane | 5.7 | ug/m ³ | 2.46 | - | U | Yes |
| 1,4-Dioxane | 18 | ug/m ³ | 2.46 | - | U | Yes |
| Bromodichloromethane | 8.2 | ug/m ³ | 2.46 | - | U | Yes |
| cis-1,3-Dichloropropene | 5.6 | ug/m ³ | 2.46 | - | U | Yes |
| 4-Methyl-2-pentanone | 7.8 | ug/m ³ | 2.46 | - | - | Yes |
| Toluene | 8.1 | ug/m ³ | 2.46 | - | - | Yes |

| | | | | | | |
|---------------------------|------|-------------------|------|---|---|-----|
| trans-1,3-Dichloropropene | 17 | ug/m ³ | 2.46 | - | - | Yes |
| 1,1,2-Trichloroethane | 3.1 | ug/m ³ | 2.46 | J | J | Yes |
| Tetrachloroethene | 8.3 | ug/m ³ | 2.46 | - | U | Yes |
| 2-Hexanone | 2.2 | ug/m ³ | 2.46 | J | J | Yes |
| Dibromochloromethane | 10 | ug/m ³ | 2.46 | - | U | Yes |
| 1,2-Dibromoethane (EDB) | 9.4 | ug/m ³ | 2.46 | - | U | Yes |
| Chlorobenzene | 5.7 | ug/m ³ | 2.46 | - | U | Yes |
| Ethyl Benzene | 2.0 | ug/m ³ | 2.46 | J | J | Yes |
| m,p-Xylene | 7.8 | ug/m ³ | 2.46 | - | - | Yes |
| o-Xylene | 3.9 | ug/m ³ | 2.46 | J | J | Yes |
| Styrene | 0.64 | ug/m ³ | 2.46 | J | J | Yes |
| Bromoform | 13 | ug/m ³ | 2.46 | - | U | Yes |
| Cumene | 1.4 | ug/m ³ | 2.46 | J | J | Yes |
| 1,1,2,2-Tetrachloroethane | 0.48 | ug/m ³ | 2.46 | J | J | Yes |
| Propylbenzene | 0.66 | ug/m ³ | 2.46 | J | J | Yes |
| 4-Ethyltoluene | 3.1 | ug/m ³ | 2.46 | J | J | Yes |
| 1,3,5-Trimethylbenzene | 0.74 | ug/m ³ | 2.46 | J | J | Yes |
| 1,2,4-Trimethylbenzene | 2.7 | ug/m ³ | 2.46 | J | J | Yes |
| 1,3-Dichlorobenzene | 15 | ug/m ³ | 2.46 | - | - | Yes |
| 1,4-Dichlorobenzene | 14 | ug/m ³ | 2.46 | - | - | Yes |
| alpha-Chlorotoluene | 6.4 | ug/m ³ | 2.46 | - | U | Yes |
| 1,2-Dichlorobenzene | 7.4 | ug/m ³ | 2.46 | - | U | Yes |
| 1,2,4-Trichlorobenzene | 36 | ug/m ³ | 2.46 | - | U | Yes |
| Hexachlorobutadiene | 52 | ug/m ³ | 2.46 | - | U | Yes |
| Naphthalene | 13 | ug/m ³ | 2.46 | - | U | Yes |

Project Number: 1701428BDate: 01/24 & 26/17

REVIEW OF VOLATILE ORGANIC PACKAGE

The following guidelines for evaluating volatile organics were created to delineate required validation actions. This document will assist the reviewer in using professional judgment to make more informed decision and in better serving the needs of the data users. The sample results were assessed according to USEPA data validation guidance documents in the following order of precedence: QC criteria from "Compendium Method TO-15. Determination of Volatile Organic Compounds (VOCs) In Air Collected In Specially-Prepared Canisters and Analyzed By Gas Chromatography/Mass Spectrometry (GC/MS), January, 1999"; USEPA Hazardous Waste Support Branch. Validating Air Samples. Volatile Organic Analysis of Ambient Air in Canisters by Method TO-15, (SOP # HW-31. Revision #6. June, 2014). The QC criteria and data validation actions listed on the data review worksheets are from the primary guidance document, unless otherwise noted.

The hardcopied (laboratory name) Eurofins data package received has been reviewed and the quality control and performance data summarized. The data review for VOCs included:

Lab. Project/SDG No.: 1701477BSample matrix: AirNo. of Samples: 4Trip blank No.: -Field blank No.: -Equipment blank No.: -Field duplicate No.: 1701428B-16A/1701428B-17A; 1701428B-18A/1701428B-19A☒ Data Completeness☒ Laboratory Control Spikes☒ Sampling Integrity/Preservation☒ Field Duplicates☒ GC/MS Tuning☒ Calibrations☒ Internal Standard Performance☒ Compound Identifications☒ Blanks☒ Compound Quantitation☒ Surrogate Recoveries☒ Quantitation LimitsOverall Comments: VOCs (full suite) by method TO-15

Definition of Qualifiers:

J- Estimated results

U- Compound not detected

R- Rejected data

UJ- Estimated nondetect

Reviewer: Rafael InfanteDate: 03/15/17

DATA COMPLETENESS

MISSING INFORMATION

DATE LAB. CONTACTED

DATE RECEIVED

A dashed diagonal line runs from the top-left corner to the bottom-right corner of the page. The line is composed of short, light gray dashes. The background of the page is white with horizontal gray lines.

All criteria were met X
 Criteria were not met
 and/or see below

SAMPLE INTEGRITY AND PRESERVATION

Canister used for sampling of the ambient air must be demonstrated clean, and leak free prior to sample collection. Cleanliness is demonstrated by the analysis of an individual canister or analysis of a representative canister, if only batch cleaning was required. Leak proof testing is performed on individual canisters. Canisters are used in conjunction with gauges, valves and flow controllers. Therefore, canister should be demonstrated clean and leak free inclusive of these components as appropriate.

- a. Leak proof test:
- Was the pressure of each canister measured before shipping? Yes or No
- Was the pressure of each canister measured before sampling? Yes or No
- Did the canister hold vacuum/pressure within +/- 2 psi from the date shipped to the sampling date? Yes or No

Note:

- The laboratory should be notified if the difference between the laboratory and field pressure is greater than 2 psi.

Actions:

Actions for use of canisters with failing leak test criteria are indicated in Table 1 below.

Table 1. Canister Leak test Actions for TO-15 Analysis*

| Matrix | Difference in initial and 24 hour pressure (psi) Criteria | Action | |
|--------|---|-------------------------------|-----------------------------------|
| | | Detected Associated Compounds | Non-Detected Associated Compounds |
| Air | ≤ 5 | No qualification | |
| Air | > 5 | J | UJ or R |

*Excessive time period (> 3months) elapsed between leak test and actual use should be considered in evaluation of canister integrity.

- b. Cleanliness
- Integrity of the canister used for sampling of air for analysis should be maintained at all times including time of shipment to the field, sampling, shipping back to the laboratory and time of analysis. Analytical results of canister cleaning verification must be taken into account in the validation of sample results.

Does the canister meet the cleanliness criteria? Yes or No

Is the canister verification included in the data package?

Yes or **No**

Actions:

Canister contamination actions are stated in Table 2 below.

Note: Laboratory stated that the SUMMA canisters employed were 100 % certified.

Table 2. Canister Contamination Actions for TO-15 Analyses

| Contamination Type/level | Canister Cleaning Result | Sample Result | Action for Samples |
|--------------------------|--------------------------|---|---|
| Clean Canister analysis | Detects | Analytes found in clean canister analysis are non-detects | No qualification required |
| | <CRQL | < CRQL | Report CRQL value with a U |
| | | \geq CRQL and < 2x the CRQL | Report concentration of sample with a U |
| | | \geq 2x the CRQL | No qualification required |
| | > CRQL | < CRQL | Report CRQL value with a U |
| | | \geq CRQL and \leq clean canister value | Report clean canister value with a U |
| | | \geq CRQL and > clean canister value | No qualification required |
| | = CRQL | \leq CRQL | Report CRQL value with a U |
| | | > CRQL | No qualification required |

c. Holding time and sample integrity

SUMMA canisters are to minimize sample charges or loss for majority of the analyte. Sample integrity is maintained by ensuring the system is closed tight and canister pressure from the time of sampling to the time of analysis is maintained within a difference allowable due to temperature change.

Was the canister pressure measured at the conclusion of the sampling period?

Yes or No

Was the canister pressure measured upon arrival to the laboratory? **Yes** or No

Was the canister pressure difference between sampling and analysis less than 5 psi? **Yes** or No

Actions:

Qualify sample results using technical holding time information as stated in Table 3.

Pressure difference between sampling and analysis should be less than 5 psi. Qualify samples as per Table 3 requirements.

Table 3. Holding Time Actions for TO-15 Volatile Analyses

| Matrix | Preserved (Pressure difference between sampling and analysis ≤ 5 psi) | Criteria | Action | |
|--------|---|--------------|-------------------------------------|--------------------------------------|
| | | | Detected Associated Compounds | Non-Detected Associated Compounds |
| Air | Yes | < 30 days | No qualification | |
| | Yes | >30 days | J | UJ |
| Air | No | < 30 days | J | UJ |
| | No | >30 days | J | R |

Complete table for all samples and note the integrity and/or preservation not within criteria

| SAMPLE ID | DATE SAMPLED | DATE ANALYZED | Pressure difference < 5 psi | ACTION |
|--|-----------------|------------------|-----------------------------------|--------|
| | | | | |
| | | | | |
| | | | | |
| All samples analyzed within the recommended method holding time. All summa canisters received in good conditions. Pressure difference < 5 psi between sampling and analysis. | | | | |
| | | | | |
| | | | | |
| | | | | |
| | | | | |
| | | | | |
| | | | | |
| | | | | |

The following pressure conversion is used, if necessary

| PRESSURE CONVERSION TABLE | | | | | | | | |
|---------------------------|-----------|---------------------|---------------------|---------|---------|---------|----------|---------------------|
| PSI | ATM | kgf/cm ² | in.H ₂ O | mmHg | in.Hg | Kpa | Bar | mm H ₂ O |
| 1 | 0.068046 | 0.070307 | 27.7276 | 51.715 | 2.03602 | 6.895 | 0.6895 | 704.28104 |
| 14.696 | 1 | 1.0332 | 407.484 | 760 | 29.921 | 101.325 | 1.01325 | 10350.0936 |
| 14.2233 | 0.96784 | 1 | 394.38 | 735.559 | 28.959 | 98.096 | 0.98067 | 10000 |
| 0.036092 | 0.002454 | 0.00253 | 1 | 1.8651 | 0.07343 | 0.249 | 0.00249 | 25.4 |
| 0.019336 | 0.001315 | 0.001359 | 0.53616 | 1 | 0.03937 | 0.1333 | 0.001333 | 13.618464 |
| 0.491154 | 0.0033421 | 0.03453 | 13.6185 | 25.4 | 1 | 3.3864 | 0.033864 | 345.9099 |
| 0.145 | 0.00987 | 0.010197 | 4.0186 | 7.5006 | 0.2953 | 1 | 0.01 | 102.07244 |
| 14.5038 | 0.98692 | 1.01972 | 402.156 | 750.062 | 29.53 | 100 | 1 | 10214.7624 |

All criteria were met X
 Criteria were not met see below

GC/MS TUNING

The assessment of the tuning results is to determine if the sample instrumentation is within the standard tuning QC limits

Gas Chromatograph/Mass Spectrometer (GC/MS) Instrument Performance Check

Action:

NOTES: This requirement does not apply when samples are analyzed by the Selected Ion Monitoring (SIM) technique.

All mass spectrometer instrument conditions must be identical to those used during the sample analysis. Background subtraction actions resulting in spectral distortions for the sole purpose of meeting the method specifications are contrary to the Quality Assurance (QA) objectives, and are therefore unacceptable.

NOTES: No data should be qualified based on BFB or DFTTP failure. Instances of this should be noted in the narrative.

All ion abundance ratios must be normalized to m/z 95, the nominal base peak, even though the ion abundance of m/z 174 may be up to 120% that of m/z 95.

1. If samples are analyzed without a preceding valid instrument performance check, qualify all data in those samples as unusable (R).
2. If the laboratory has made minor transcription errors which do not significantly affect the data, the data reviewer should make the necessary corrections on a copy of the form.
3. If the laboratory has failed to provide the correct forms or has made significant transcription or calculation errors, the Region's designated representative should contact the laboratory and request corrected data. If the information is not available, the reviewer must use professional judgment to assess the data and notify the Project Officer (PO).
4. If ion abundance criteria are not met, professional judgment may be applied to determine to what extent the data may be utilized. When applying professional judgment to this topic, the most important factors to consider are the empirical results that are relatively insensitive to location on the chromatographic profile and the type of instrumentation. Therefore, the critical ion abundance criteria for BFB are the m/z 95/96, 174/175, 174/176, and 176/177 ratios. The relative abundances of m/z 50 and 75 are of lower importance. This issue is more critical for Tentatively Identified Compounds (TICs) than for target analytes.
5. Note, in the Data Review Narrative, decisions to use analytical data associated with BFB instrument performance check failures (not meeting contract requirements).
6. If the reviewer has reason to believe that instrument performance check criteria were achieved using techniques other than those described in the Compendium method TO-15 entitled "Determination Of Volatile Organic Compounds(VOCs) In Air Collected In Specially-Prepared Canisters And Analyzed By Gas Chromatography/Mass Spectrometry(GC/MS)", section 10.4, obtain additional information on the instrument performance checks. If the techniques employed are found to be at variance with the contract requirements, the performance and procedures of the laboratory may merit evaluation.
7. Use professional judgment to determine whether associated data should be qualified based on the spectrum of the mass calibration compound.

☒ The BFB performance results were reviewed and found to be within the specified criteria.

☒ BFB tuning was performed for every 24 hours of sample analysis.

If no, use professional judgment to determine whether the associated data should be accepted, qualified or rejected.

List the samples affected:

If mass calibration is in error, all associated data are rejected.

All criteria were met X
 Criteria were not met
 and/or see below

CALIBRATION VERIFICATION

Compliance requirements for satisfactory instrument calibration are established to ensure that the instrument is capable of producing and maintaining acceptable quantitative data.

Date of initial calibration: 12/01/2016
 Date of initial calibration verification: -
 Dates of continuing calibration: 02/02/17(10:10_AM);_02/06/17_(9:15_AM)
 Instrument ID numbers: MSD-14
 Matrix/Level: Air/low

| DATE | LAB FILE ID# | CRITERIA OUT RFs, %RSD, %D, r | COMPOUND | SAMPLES AFFECTED |
|--|--------------|----------------------------------|-----------|-------------------------------|
| | | | | |
| Initial and continuing calibrations meet method specific requirements except for the cases described in this document. | | | | |
| 02/06/17 | 1701428B-21B | 35 % | Freon 11 | 1701428B-16A; 1701428B-17A |
| | | 34 % | Freon 113 | |
| | | | | |
| | | | | |

Note: Results qualified (J or UJ) in affected sample.

The following criteria apply:

Table 5. Initial Calibration Actions for TO-15 Analyses

| Criteria for TO-15 Analysis | Action | |
|---|---|---|
| | Detected Associated Compounds | Non-Detected Associated Compounds |
| RRF < 0.010 (poor response volatile target compounds, Table 4) RRF < 0.050 (all other volatile target compounds) | J (based on mass spectral identification) | R |
| RRF > 0.010 (poor response volatile target compounds, Table 4) RRF > 0.050 (all other volatile target compounds) | No qualification | |
| % RSD > 40.0 or < -40.0 (poor response volatile target compounds, Table 4) % RSD > 30.0 or < -30.0 (all other Volatile target compounds) | No qualification | |
| % RSD < 40.0 and > -40.0 (poor response volatile target compounds, Table 4) % RSD < 30.0 and > -30.0 (all other volatile target compounds) | J | Use professional judgment |

Table 6. Continuing Calibration Verification (CCV) Actions for TO-15 Analyses

| Criteria for CCV | Action | |
|---|---|-----------------------------------|
| | Detected Associated Compounds | Non-Detected Associated Compounds |
| RRF < 0.010 (poor response volatile target compounds, Table 4) RRF < 0.050 (all other volatile target compounds) | J (based on mass spectral identification) | R |
| RRF > 0.010 (poor response volatile target compounds, Table 4) RRF > 0.050 (all other volatile target compounds) | No qualification | |
| %D > 40.0 or < -40.0 (poor response volatile target compounds, Table 4) %D > 30.0 or < -30.0 (all other Volatile target compounds) | J | UJ |
| %D < 40.0 and > -40.0 (poor response volatile target compounds, Table 4) %D < 30.0 and > -30.0 (all other volatile target compounds) | No qualification | |

If the % D for daily calibration exceeds -90, use professional judgment to see if non-detects need to be qualified as unusable "R"

A separate worksheet should be filled for each initial curve

Table 4. TO 15 Volatile Compounds List*

| Compound | CAS Number | Synonyms |
|-------------------------------|-----------------|---|
| Acetone | 67-64-1 | Dimethyl ketone; Dimethylformaldehyde; 2-Propanone |
| Allyl chloride | 107-05-1 | 3-Chloropropene; 3-Chloroprene |
| Benzene | 71-43-2 | Benzol; Benzine |
| Benzyl chloride | 100-44-7 | Chloromethylbenzene; alpha-Chlorotoluene |
| Bromodichloromethane | 75-27-4 | Monobromodichloromethane; Methane-bromodichloro |
| Bromoethene | 593-60-2 | Vinyl bromide; Monobromoethene |
| Bromoform | 75-25-2 | Tribromoethane |
| Bromomethane | 74-83-9 | Methyl bromide; Monobromomethane |
| 1,3-Butadiene | 106-99-0 | Biethylene; Erythrene; Pyrrolyene |
| Carbon disulfide | 75-15-0 | Carbon bisulfide; Carbon sulfide |
| Carbon tetrachloride | 56-23-5 | Carbon tet; Tetrachloromethane |
| Chlorobenzene | 108-90-7 | Monochlorobenzene; Chlorobenzol; Benzene chloride |
| Chloroethane | 75-00-3 | Ethyl chloride; Chlorene; Chloryl |
| Chloroethene | 75-01-4 | Vinyl chloride; Ethylene monochloride |
| Chloroform | 67-66-3 | Trichloromethane; Methyltrichloride; Methane trichloride |
| Chloromethane | 74-87-3 | R40; Methyl chloride; Monochloromethane |
| Cyclohexane | 110-82-7 | Hexamethylene; Hexahydrobenzene; Hexanaphthene |
| Dibromochloromethane | 124-48-1 | Chlorodibromomethane |
| 1,2-Dibromoethane | 106-93-4 | EDB; Ethylene dibromide |
| 1,2-Dichlorobenzene | 95-50-1 | ODB; Chloroben |
| 1,3-Dichlorobenzene | 541-73-1 | meta-Dichlorobenzene; m-Phenylenedichloride |
| 1,4-Dichlorobenzene | 106-46-7 | para-Dichlorobenzene; Parazene; Santochlor |
| 1,1-Dichloroethane | 75-34-3 | Ethylidene chloride; Ethylidene dichloride |
| 1,2-Dichloroethane | 107-06-2 | Ethylene dichloride; Glycol dichloride; 1,2-DCA |
| 1,1-Dichloroethene | 75-35-4 | 1,1-DCE; Vinylidene chloride |
| cis-1,2-Dichloroethylene | 156-59-2 | cis-1,2-DCE; cis-Acetylene dichloride |
| trans-1,2-Dichloroethylene | 156-60-5 | trans-1,2-DCE; trans-Acetylene dichloride |
| 1,2-Dichloropropane | 78-87-5 | Propylene dichloride; Propylene chloride |
| cis-1,3-Dichloropropene | 10061-01-5 | 1-Propene, 1,3-dichloro-, (z)-; cis-1,3-Dichloro-1-Propene |
| trans-1,3-Dichloropropene | 10061-02-6 | trans-1,3-Dichloro-1-Propene; trans-1,3-Dichloropropylene |
| 1,4-Dioxane | 123-91-1 | Diethylene dioxide; Diethylene ether |
| Ethyl acetate | 141-78-6 | Acetic acid ethyl ester; Acetic ether |
| Ethylbenzene | 100-41-4 | Ethylbenzol; Phenylethane |
| 4-Ethyltoluene | 622-96-8 | 1-Ethyl-4-methyl benzene; p-Methylethylbenzene |
| Freon 11 (CCl ₃ F) | 75-69-4 | Trichlorofluoromethane; Fluorotrichloromethane; Fluorocarbon 11 |

| | | |
|--|----------------|--|
| Freon 12 (CCl ₂ F ₂) | 75-71-8 | Dichlorodifluoromethane; Fluorocarbon 12 |
| Freon 113 (C ₂ Cl ₃ F ₃) | 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane; Fluorocarbon 113; 1,1,2-Trichlorotrifluoroethane |
| Freon 114 (C ₂ Cl ₂ F ₄) | 76-14-2 | 1,2-Dichlorotetrafluoroethane; Halocarbon 114; 1,2-Dichloro-1,1,2,2-tetrafluoroethane |
| Heptane | 142-82-5 | Dipropylmethane; Heptyl hydride |
| Hexachlorobutadiene | 87-68-3 | 1,3-Hexachlorobutadiene; Perchlorobutadiene |
| Hexane | 110-54-3 | n-Hexane; Hexyl hydride |
| 2-Hexanone | 591-78-6 | Methyl butyl ketone; Butyl methyl ketone; Hexan-2-one |
| Isopropyl alcohol | 67-63-0 | 2-Propanol; Isopropanol |
| Methylene chloride | 75-09-2 | Dichloromethane; Methylene dichloride |
| Methyl ethyl ketone | 78-93-3 | MEK; 2-Butanone; Ethyl methyl ketone |
| Methyl isobutyl ketone | 108-10-1 | MIBK; 2-Pentanone; Hexone; Isopropylacetone |
| Methyl tert-butyl ether | 1634-04-4 | MTBE; 2-Methoxy-2-methylpropane; tert-Butyl methyl ether |
| Propylene | 115-07-1 | Propene; Methylene |
| Styrene | 100-42-5 | Vinylbenzene; Phenylethylene |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | Tetrachloroethane; Acetylene tetrachloride; Bonoform |
| Tetrachloroethene | 127-18-4 | PCE; PERC; Perchloroethylene; Ethylene tetrachloride; Carbon bichloride; Carbon dichloride |
| Tetrahydrofuran | 109-99-9 | Diethylene oxide; Butylene oxide |
| Toluene | 108-88-3 | Toluol; Methylbenzene |
| 1,2,4-Trichlorobenzene | 120-82-1 | 1,2,4-Trichlorobenzol |
| 1,1,1-Trichloroethane | 71-55-6 | Methyl chloroform; Trichloroethane |
| 1,1,2-Trichloroethane | 79-00-5 | beta-Trichloroethane; Ethane trichloride; Vinyl trichloride |
| Trichloroethene | 79-01-6 | TCE; Acetylene trichloride; Ethinyl trichloride |
| 1,2,4-Trimethylbenzene | 95-63-6 | Pseudocumene; Pseudocumol |
| 1,3,5-Trimethylbenzene | 108-67-8 | Mesitylene; Trimethylbenzol |
| 2,2,4-Trimethylpentane | 540-84-1 | Iso-octane; Isobutyltrimethylmethane |
| Vinyl acetate | 108-05-4 | Acetic acid ethenyl ether; Ethenyl acetate |
| p-Xylene | 106-42-3 | p-Methyltoluene; 1,4-dimethylbenzene |
| m-Xylene | 108-38-3 | m-Methyltoluene; 1,3-dimethylbenzene |
| o-Xylene | 95-47-6 | o-Methyltoluene; 1,2-Dimethylbenzene |

*Laboratories use different sets and subsets of analytes on as needed basis.

NOTES:

Compounds in bold italicized letters may have poor GCMS response. These poor response compounds are evaluated using more relaxed relative response factor criteria as stated below.

All criteria were met _____
 Criteria were not met _____
 and/or see below _____X_____

V A. BLANK ANALYSIS RESULTS (Sections 1 & 2)

The assessment of the blank analysis results is to determine the existence and magnitude of contamination problems. The criteria for evaluation of blanks apply only to blanks associated with the samples, including trip, equipment, and laboratory blanks. If problems with any blanks exist, all data associated with the case must be carefully evaluated to determine whether or not there is an inherent variability in the data for the case, or if the problem is an isolated occurrence not affecting other data.

List the contamination in the blanks below. High and low levels blanks must be treated separately.

Blanks criteria and appropriate actions

Table 7. Blank Actions for TO-15 Analyses

| Blank Type | Blank Result | Sample Result | Action for Samples |
|---|------------------------|---|--|
| Method, Storage, Field, Trip, Instrument*** | Detects | Not detected | No qualification required |
| | < CRQL * | < CRQL* | Report CRQL value with a U |
| | | \geq CRQL* and < 2x the CRQL** | Report concentration of sample with a "U" |
| | | \geq 2x the CRQL** | No qualification required |
| | > CRQL * | < CRQL * | Report CRQL value with a U |
| | | \geq CRQL* and \leq blank concentration | Report blank value for sample concentration with a U |
| | | \geq CRQL* and > blank concentration | No qualification required |
| | = CRQL* | \leq CRQL * | Report CRQL value with a U |
| | | > CRQL * | No qualification required |
| | Gross contamination ** | Detects | Report blank value for sample concentration with a U |

* 2x the CRQL for methylene chloride, 2-butanone and acetone.

** 4x the CRQL for methylene chloride, 2-butanone, and acetone.

*** Qualifications based on instrument blank results affect only the sample analyzed immediately after the sample that has target compounds that exceed the calibration range or non-target compounds that exceed 100 µg/L.

| DATE ANALYZED | LAB ID | LEVEL/ MATRIX | COMPOUND | CONCENTRATION UNITS |
|---|--------------|---------------|----------|---------------------|
| All method blank meet method specific criteria except for the cases described in this document. | | | | |
| 02/03/17 | 1701428B-20A | Air/low | Acetone | 0.58 ppbv |
| | | | Benzene | 0.072 ppbv |

Note: Concentration detected below the reporting limit, results qualified following the table above.

Field/Equipment/Trip blank

| DATE ANALYZED | LAB ID | LEVEL/MATRIX | COMPOUND | CONCENTRATION UNITS |
|---------------|--------|--------------|----------|---------------------|
|---------------|--------|--------------|----------|---------------------|

No field/trip/equipment blanks analyzed with this data package.

Field/trip blank actions

Field or Trip blank when available should be assessed for possible contaminants in the canister used for trip blank. This canister and its analytical results are specific to the trip blank sample **only**. If contaminants are present in the canister used for trip blank, its suitability for use as trip blank can be assessed using the following criteria.

Table 8. Field/Trip Blank suitability based on Canister contamination

| Clean canister Result | Field/Trip Blank Result | Action for Field/Trip Blank |
|-----------------------|--|---|
| Detects | Not detected | No qualification, no action for samples is required |
| Detects | < clean canister result or \geq clean canister result but < 2X the clean canister result | Report as non-detect "U", invalid as trip blank, no action for samples is required. |
| | $\geq 2x$ the clean canister result | No qualification, valid trip blank for sample actions. |

| CONTAMINATION SOURCE/LEVEL | COMPOUND | CONC/UNITS | AL/UNITS | SQL | AFFECTED SAMPLES |
|----------------------------|----------|------------|----------|-----|------------------|
| | | | | | |
| | | | | | |
| | | | | | |
| | | | | | |

All criteria were met X
 Criteria were not met
 and/or see below

SURROGATE SPIKE RECOVERIES

Laboratory performance of individual samples is established by evaluation of surrogate spike recoveries. All samples are spiked with surrogate compounds prior to sample analysis. The accuracy of the analysis is measured by the surrogate percent recovery. Since the effects of the sample matrix are frequently outside the control of the laboratory and may present relatively unique problems, the validation of data is frequently subjective and demands analytical experience and professional judgment.

List the percent recoveries (%Rs) which do not meet the criteria for surrogate recovery.

Matrix: solid/aqueous

| SAMPLE ID | SURROGATE COMPOUND | | | ACTION |
|-----------|-----------------------|------------|-------|--------|
| | 1,2-DICHLOROETHANE-d4 | Toluene-d8 | 4-BFB | |

 Surrogate recoveries within laboratory control limits

QC Limits* (Air)

 LL to UL 70 to 130 70 to 130 70 to 130

- * QC limits are laboratory in-house performance criteria, LL = lower limit, UL = upper limit.
- * If QC limits are not available, use limits of 80 – 120 % for aqueous and 70 – 130 % for solid samples.

Actions:

| QUALITY | %R < 10% | %R = 10% - LL | %R > UL |
|--------------------|----------|---------------|---------|
| Positive results | J | J | J |
| Nondetects results | R | UJ | Accept |

Surrogate action should be applied:

If one or more surrogate in the VOC fraction is out of specification, but has a recovery of > 10%.

If any one surrogate in a fraction shows < 10 % recovery.

All criteria were met X
 Criteria were not met
 and/or see below

VIII. LABORATORY CONTROL SAMPLE (LCS) ANALYSIS

This data is generated to determine accuracy of the analytical method for various matrices. LCS concentration should be in the middle of the calibration range and under the same sample conditions.

1. LCS Recoveries Criteria

Table 9. LCS/LCSD Actions for TO-15 Analyses

| Criteria | Action | |
|--|-------------------------------|-----------------------------------|
| | Detected Associated Compounds | Non-detected Associated Compounds |
| Percent recovery Criteria | | |
| %R > Upper Acceptance Limit (>130%) | J | No qualification |
| %R in the acceptable range, 70-130% | No qualification | |
| %R < Lower Acceptance Limit (< 70 %) | J | UJ |
| %R < 50% | J | R |
| Lower Acceptance Limit ≤ %R ≤ Upper Acceptance Limit | No qualification | |
| | | |
| Relative Percent Difference Criteria | | |
| % RPD ≤ 25% | No qualification | |
| % RPD > 25 % | J | UJ |

| LCS ID | COMPOUND | % R | QC LIMIT |
|---|-------------|-------|------------|
| LCS/LCSD (Blank spike) analyzed in this data package. % recoveries and RPD _____ | | | |
| within laboratory control limits except for the cases described in this document. _____ | | | |
| 1701428B-22B | Freon 113 | 131 % | 70 - 130 % |
| 1701428B-22BB | Naphthalene | 152 % | 60 - 140 % |
| | | | |
| | | | |
| | | | |
| | | | |

Note: Detected compounds are qualified with a J qualifier in affected samples.

2. Frequency Criteria:

Where LCS analyzed at the required frequency and for each matrix? **Yes** or No.

If no, the data may be affected. Use professional judgment to determine the severity of the effect and qualify data accordingly. Discuss any actions below and list the samples affected.

All criteria were met _____
 Criteria were not met _____
 and/or see below _____X_____

IX. FIELD/LABORATORY DUPLICATE PRECISION

Sample IDs: _____ LCS/LCSD (lab) _____

Matrix: _____ Air _____

Sample IDs: _____ 1701428B-16A/1701428B-17A _____

Matrix: _____ Air _____

Sample IDs: _____ 1701428B-18A/1701428B-19A _____

Matrix: _____ Air _____

Field duplicate samples may be taken and analyzed as an indication of overall precision. These analyses measure both field and lab precision; therefore, the results may have more variability than laboratory duplicates which only laboratory performance. It is also expected that soil duplicate results will have a greater variance than water matrices due to difficulties associated with collecting identical field duplicate samples.

The project QAPP should be reviewed for project-specific information. In the absence of QAPP guidance for validated data from field duplicates, the following action will be taken.

Identify which samples within the data package are field duplicates. Estimate the relative percent difference (RPD) between the values for each compound. Note large RPDs (>50 %) in the narrative. Use professional judgment to qualify data when RPD > 50 %.

| COMPOUND | SQL | SAMPLE CONC. | DUPLICATE CONC. | RPD | ACTION |
|--|-----|--------------|-----------------|-----|--|
| 1701428B-18A/1701428B-19A | | | | | |
| Ethanol | 9.4 | 270 | 460 | 52 | Qualify results (J in affected samples |
| 2-propanol | 12 | 460 | 880 | 63 | |
| Laboratory/field duplicate analyzed with this data package. RPD within laboratory and method performed criteria for analytes with concentration > 5 x SQL except for the cases described in this document. | | | | | |

Other suggested actions:

Qualify as estimated positive results (J) and nondetects (UJ) for the compound that exceeded the above criteria. For organics, only the sample and duplicate will be qualified.

If an RPD cannot be calculated because one or both of the sample results is not detected, the following actions apply:

If one sample result is not detected and the other is greater than 5x the SQL qualify (J/UJ).

If one sample value is not detected and the other is greater than 5x the SQL and the SQLs for the sample and duplicate are significantly different, use professional judgment to determine if qualification is appropriate.

If one sample value is not detected and the other is less than 5x, use professional judgment to determine if qualification is appropriate.

If both sample and duplicate results are not detected, no action is needed.

All criteria were met X
 Criteria were not met
 and/or see below

X. INTERNAL STANDARD PERFORMANCE

The assessment of the internal standard (IS) parameter is used to assist the data reviewer in determining the condition of the analytical instrumentation.

List the internal standard area of samples which do not meet the criteria.

- * Area of +40% or -40% of the IS area in the associated calibration standard (CCV standard or mid-point from initial calibration).
- * Retention time (RT) within ± 20 seconds of the IS area in the associated calibration standard.

Table 10. Internal Standard Actions for TO-15 Analyses

| Criteria | Action | |
|--|--------------------------------|------------------------------------|
| | Detected Associated Compounds* | Non-detected Associated Compounds* |
| Area counts > 140% of CCV or mid-point standard from initial calibration) | J- | No qualification |
| Area counts < 60% of CCV or mid-point standard from initial calibration) | J+ | R |
| Area counts $\geq 60\%$ but $\leq 140\%$ of CCV or mid-point standard from initial calibration) | No qualification | |
| RT difference > 20.0 seconds between samples CCV or mid-point standard from initial calibration) | R* | |
| RT difference < 20.0 seconds between samples and CCV or mid-point standard from initial calibration) | No qualification | |

* Examine the chromatographic profile for that sample to determine if any false positives or negatives exist. For shifts of a large magnitude, the reviewer may consider partial or total rejection of the data for that sample fraction. Detects should not need to be qualified as unusable (R) if the mass spectral criteria are met.

| | | | | | |
|------|-----------|--------|---------|------------------|--------|
| DATE | SAMPLE ID | IS OUT | IS AREA | ACCEPTABLE RANGE | ACTION |
|------|-----------|--------|---------|------------------|--------|

Internal standard area and retention times within laboratory control limits for both samples and calibration standards

All criteria were met ☒X____
Criteria were not met
and/or see below _____

XII. SAMPLE QUANTITATION

The sample quantitation evaluation is to verify laboratory quantitation results. In the space below, please show a minimum of one sample calculation:

1701428B-16A

Methyl-tert-butyl-ether

RF = 3.20626

$$[] = (7212)(400)/(87894)(3.20626)$$

$$= 10.24 \text{ ppbv OK}$$

All criteria were met ☒X___
 Criteria were not met
 and/or see below _____

XII. QUANTITATION LIMITS

A. Dilution performed

| SAMPLE ID | DILUTION FACTOR | REASONS FOR DILUTION |
|--|-----------------|----------------------|
| Samples diluted by a factor of 2.50 or less except for the followings. | | |
| 1701428B-16A | 24.9 | Matrix interference |
| 1701428B-17A | 24.2 | Matrix interference |
| | | |
| | | |
| | | |
| | | |
| | | |
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| | | |

System Performance

Action:

Use professional judgment to qualify the data if it is determined that system performance has degraded during sample analyses. Note, for Laboratory Project Officer (PO) action, any degradation of system performance which significantly affected the data.

Note:

Overall Assessment of Data

Action:

1. Use professional judgment to determine if there is any need to qualify data which were not qualified based on the Quality Control (QC) criteria previously discussed.
2. Write a brief narrative to give the user an indication of the analytical limitations of the data. Note, for Laboratory Project Officer (PO) action, any inconsistency of the data with the Sample Delivery Group (SDG) Narrative. If sufficient information on the intended use and required quality of the data is available, the reviewer should include their assessment of the usability of the data within the given context. This may be used as part of a formal Data Quality Assessment (DQA).

Overall assessment of the data: Results are valid; the data can be used for decision making purposes.